

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061319\
 Data File : BG041224.D
 Acq On : 13 Jun 2019 10:13
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:48:10 AM

Quant Time: Jun 13 10:51:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 23:08:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	54988	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	236294	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	162314	20.00	ng	0.00
64) Phenanthrene-d10	17.49	188	397146	20.00	ng	0.00
76) Chrysene-d12	21.78	240	401330	20.00	ng	0.02
87) Perylene-d12	25.14	264	419603	20.00	ng	0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	239250	78.46	ng	0.00
7) Phenol-d6	7.25	99	366838	77.89	ng	0.00
23) Nitrobenzene-d5	9.29	82	428665	80.54	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	195619	81.18	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	947934	84.10	ng	0.00
79) Terphenyl-d14	20.09	244	1587575	83.96	ng	0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	49725	35.934	ng	98
3) Pyridine	3.91	79	144099	35.193	ng	98
4) n-Nitrosodimethylamine	3.83	42	82189	43.250	ng	# 78
6) Aniline	7.43	93	244786	38.940	ng	97
8) 2-Chlorophenol	7.66	128	143296	39.416	ng	95
9) Benzaldehyde	7.24	77	116527	41.478	ng	90
10) Phenol	7.28	94	195717	38.759	ng	99
11) bis(2-Chloroethyl)ether	7.52	93	142751	38.969	ng	97
12) 1,3-Dichlorobenzene	7.99	146	175390	40.393	ng	96
13) 1,4-Dichlorobenzene	8.14	146	179923	40.936	ng	94
14) 1,2-Dichlorobenzene	8.46	146	169224	39.654	ng	94
15) Benzyl Alcohol	8.35	79	169844	38.986	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.62	45	226359	37.815	ng	100
17) 2-Methylphenol	8.54	107	140425	38.408	ng	97
18) Hexachloroethane	9.19	117	70316	41.509	ng	97
19) n-Nitroso-di-n-propylamine	8.91	70	138812	38.301	ng	98
20) 3+4-Methylphenols	8.87	107	190922	37.699	ng	98
22) Acetophenone	8.94	105	264213	39.860	ng	98
24) Nitrobenzene	9.33	77	220344	40.622	ng	96
25) Isophorone	9.84	82	386560	38.162	ng	99
26) 2-Nitrophenol	10.04	139	95904	42.888	ng	90
27) 2,4-Dimethylphenol	10.08	122	133540	36.159	ng	92
28) bis(2-Chloroethoxy)methane	10.33	93	209046	38.237	ng	98
29) 2,4-Dichlorophenol	10.57	162	177141	37.771	ng	93
30) 1,2,4-Trichlorobenzene	10.78	180	219083	41.064	ng	97
31) Naphthalene	10.98	128	505056	39.809	ng	99
32) Benzoic acid	10.25	122	90118	34.196	ng	92
33) 4-Chloroaniline	11.10	127	223841	38.026	ng	96
34) Hexachlorobutadiene	11.24	225	172014	42.138	ng	97
35) Caprolactam	11.91	113	53874	34.786	ng	89
36) 4-Chloro-3-methylphenol	12.20	107	198115	36.988	ng	93
37) 2-Methylnaphthalene	12.58	142	371179	37.987	ng	98
38) 1-Methylnaphthalene	12.58	142	371179	40.312	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	277235	42.377	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.90	237	134097	42.441	ng	95
43) 2,4,6-Trichlorophenol	13.17	196	177479	41.936	ng	97
44) 2,4,5-Trichlorophenol	13.25	196	178573	40.008	ng	99
46) 1,1'-Biphenyl	13.57	154	495853	41.270	ng	98
47) 2-Chloronaphthalene	13.63	162	435098	42.183	ng	99
48) 2-Nitroaniline	13.83	65	153497	41.482	ng	92
49) Acenaphthylene	14.47	152	627819	40.442	ng	99
50) Dimethylphthalate	14.19	163	547847	39.873	ng	98
51) 2,6-Dinitrotoluene	14.33	165	121829	41.124	ng	97
52) Acenaphthene	14.81	154	373946	39.763	ng	96
53) 3-Nitroaniline	14.66	138	118960	40.157	ng	99
54) 2,4-Dinitrophenol	14.86	184	50677	42.963	ng	94
55) Dibenzofuran	15.14	168	636468	40.221	ng	99
56) 4-Nitrophenol	14.96	139	76536	37.667	ng	92
57) 2,4-Dinitrotoluene	15.11	165	171515	40.986	ng	# 98
58) Fluorene	15.79	166	495203	39.295	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.36	232	178219	40.630	ng	96
60) Diethylphthalate	15.54	149	553900	39.502	ng	99
61) 4-Chlorophenyl-phenylether	15.77	204	330162	40.307	ng	97
62) 4-Nitroaniline	15.82	138	129714	41.361	ng	93
63) Azobenzene	16.07	77	488539	38.333	ng	98
65) 4,6-Dinitro-2-methylphenol	15.87	198	92070	44.111	ng	93
66) n-Nitrosodiphenylamine	15.99	169	442126	39.602	ng	99
67) 4-Bromophenyl-phenylether	16.67	248	212584	39.664	ng	98
68) Hexachlorobenzene	16.78	284	229099	40.142	ng	97
69) Atrazine	16.94	200	168686	39.158	ng	97
70) Pentachlorophenol	17.14	266	112033	38.951	ng	94
71) Phenanthrene	17.53	178	840890	40.649	ng	98
72) Anthracene	17.62	178	839382	41.141	ng	99
73) Carbazole	17.90	167	727736	41.570	ng	99
74) Di-n-butylphthalate	18.43	149	892448	41.017	ng	98
75) Fluoranthene	19.54	202	1086852	42.536	ng	99
77) Benzidine	19.72	184	308789	32.253	ng	96
78) Pyrene	19.90	202	1077468	41.300	ng	100
80) Butylbenzylphthalate	20.77	149	424206	41.782	ng	99
81) Benzo(a)anthracene	21.76	228	1087969	40.725	ng	98
82) 3,3'-Dichlorobenzidine	21.68	252	393046	41.830	ng	99
83) Chrysene	21.84	228	1043343	40.811	ng	98
84) Bis(2-ethylhexyl)phthalate	21.62	149	582082	40.727	ng	99
85) Di-n-octyl phthalate	22.88	149	975735	40.992	ng	100
86) Indeno(1,2,3-cd)pyrene	29.01	276	1024355m	32.709	ng	
88) Benzo(b)fluoranthene	24.07	252	1008534	39.627	ng	97
89) Benzo(k)fluoranthene	24.14	252	1091884	43.355	ng	98
90) Benzo(a)pyrene	24.98	252	971774	40.021	ng	96
91) Dibenzo(a,h)anthracene	29.07	278	838504m	34.560	ng	
92) Benzo(g,h,i)perylene	30.23	276	808413m	34.296	ng	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

